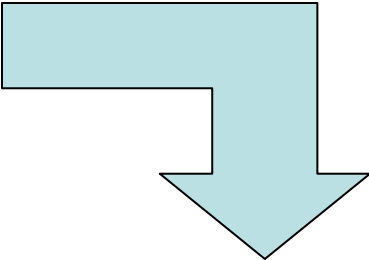


**Table 3. Individual Values of the Massic Energy of Combustion,  $\Delta_c u^\circ$ , for the Compounds, at  $T = 298.15\text{ K}$**

1-methyl-3-piperidinol	1-ethyl-3-piperidinol	1-methyl-4-piperidinol
	$-\Delta_c u^\circ/(\text{J}\cdot\text{g}^{-1})$	
34082.67	35385.57	33998.83
34086.76	35392.81	34000.16
34072.03	35393.47	34005.49
34066.73	35378.82	33978.33
34064.55	35366.63	33984.60
34060.09	35372.96	33990.18
	35388.60	
	$-\langle\Delta_c u^\circ\rangle/(\text{J}\cdot\text{g}^{-1})$	
$34072.1 \pm 4.3$	$35382.7 \pm 3.9$	$33992.9 \pm 4.2$



**Data Summary entry for this table**

<b>Table #:</b> 3
<b>System type (Pure, Binary, Ternary, Reaction):</b> Reaction
<b>Chemical Reaction(s):</b> Combustion of the liquids (See eq. 1 & 2): 1-methyl-3-piperidinol 1-ethyl-3-piperidinol 1-methyl-4-piperidinol
<b>Property:</b> Energy of Combustion
<b>Experimental Method:</b> static oxygen bomb calorimetry
<b>Combined Expanded Uncertainty (<math>k = 2</math>) for the Property:</b> $2\sigma(u)$ values are given in the table
<b>State Variables and Constraints:</b> temperature $T = 298.15\text{ K}$ , pressure $p = 100\text{ MPa}$
<b>Standard Uncertainty (<math>k = 1</math>) for Variables and Constraints:</b> $\sigma(T) = 0.01\text{ K}$ ; $\sigma(p) = 1\%$

**Note:** There is no need to mention derived properties, such as enthalpies of formation